

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092617\
 Data File : BF098836.D
 Acq On : 26 Sep 2017 17:03
 Operator : SJ/JU
 Sample : I5363-09 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-3-E-COMP

Manual Integrations
 APPROVED

Sohil
 9/27/2017 12:44:55 PM

Quant Time: Sep 27 01:55:02 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.91	152	144277	20.00	ng	-0.01
21) Naphthalene-d8	8.20	136	598727	20.00	ng	0.00
38) Acenaphthene-d10	9.95	164	269803	20.00	ng	-0.01
63) Phenanthrene-d10	11.44	188	438543	20.00	ng	0.00
75) Chrysene-d12	14.07	240	283794	20.00	ng	-0.01
86) Perylene-d12	15.57	264	271081	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	136337	15.04	ng	-0.02
7) Phenol-d6	6.53	99	168293	15.05	ng	-0.02
23) Nitrobenzene-d5	7.47	82	95875	10.27	ng	-0.02
41) 2,4,6-Tribromophenol	10.73	330	29166	13.21	ng	-0.02
44) 2-Fluorobiphenyl	9.27	172	200668	12.42	ng	-0.01
78) Terphenyl-d14	13.02	244	145755	11.68	ng	-0.02
Target Compounds						Qvalue
70) Phenanthrene	11.46	178	55948	2.383	ng	98
74) Fluoranthene	12.65	202	80743	3.397	ng	97
77) Pyrene	12.88	202	69738	3.223	ng	98
80) Benzo(a)anthracene	14.07	228	36062	2.099	ng	93
82) Chrysene	14.10	228	38248	2.338	ng	98
87) Benzo(b)fluoranthene	15.13	252	50977m	3.059	ng	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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